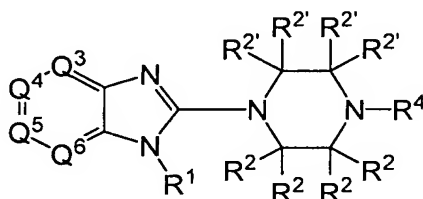


Amendments to the Claims

**This listing of claims will replace all prior versions and listings of claims in the application:**

Claim 1 (currently amended): A compound having the structure:



or any pharmaceutically-acceptable salt thereof, wherein:

m is independently at each instance 0, 1 or 2;

Q<sup>3</sup> is N or C(R<sup>5</sup>);

Q<sup>4</sup> is ~~N or~~ C(R<sup>6</sup>);

Q<sup>5</sup> is ~~N or~~ C(R<sup>6'</sup>);

Q<sup>6</sup> is N or C(R<sup>5'</sup>); wherein no more than one of Q<sup>3</sup> and Q<sup>6</sup> is N;

R<sup>1</sup> is H or -(C(R<sup>2</sup>)(R<sup>2</sup>))<sub>m</sub>-R<sup>g</sup>;

R<sup>2</sup> is, independently, in each instance, H, C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, -O(C<sub>1-7</sub>alkyl), -N(C<sub>1-7</sub>alkyl)R<sup>a</sup>, or a C<sub>1-6</sub>alkyl substituted by 1, 2 or 3 substituents selected from halo, cyano, -OR<sup>a</sup>, -OC(=O)R<sup>b</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup> and -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup>; wherein any two geminal R<sup>2</sup> groups may additionally be oxo;

R<sup>2'</sup> is, independently, in each instance, H, C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, -O(C<sub>1-7</sub>alkyl), -N(C<sub>1-7</sub>alkyl)R<sup>a</sup>, or a C<sub>1-6</sub>alkyl substituted by 1, 2 or 3 substituents selected from halo, cyano, -OR<sup>a</sup>, -OC(=O)R<sup>b</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup> and -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup>; wherein any two geminal R<sup>2'</sup> groups may additionally be oxo;

R<sup>4</sup> is R<sup>c</sup> substituted by 0, 1, 2, 3 or 4 substituents selected from R<sup>c</sup>, R<sup>e</sup>, halo, C<sub>1-4</sub>haloalkyl, cyano, nitro, -C(=O)R<sup>e</sup>, -C(=O)OR<sup>h</sup>, -C(=O)NR<sup>a</sup>R<sup>h</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>h</sup>, -OR<sup>h</sup>, -OC(=O)R<sup>e</sup>, -OC(=O)NR<sup>a</sup>R<sup>h</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>e</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>h</sup>,

-OC<sub>2-6</sub>alkylOR<sup>h</sup>, -SR<sup>c</sup>, -S(=O)R<sup>c</sup>, -S(=O)<sub>2</sub>R<sup>c</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>c</sup>,  
-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>h</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>R<sup>h</sup>, -N(R<sup>a</sup>)C(=O)R<sup>c</sup>,  
-N(R<sup>a</sup>)C(=O)OR<sup>h</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>h</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>c</sup>,  
-N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>h</sup>, -C(=O)R<sup>g</sup>, -C(=O)OR<sup>g</sup>,  
-C(=O)NR<sup>a</sup>R<sup>g</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>g</sup>, -OR<sup>g</sup>, -OC(=O)R<sup>g</sup>, -OC(=O)NR<sup>a</sup>R<sup>g</sup>,  
-OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>g</sup>, -OC(=O)N(R<sup>g</sup>)S(=O)<sub>2</sub>R<sup>c</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>g</sup>,  
-OC<sub>2-6</sub>alkylOR<sup>g</sup>, -SR<sup>g</sup>, -S(=O)R<sup>g</sup>, -S(=O)<sub>2</sub>R<sup>g</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>g</sup>, -S(=O)<sub>2</sub>N(R<sup>g</sup>)C(=O)R<sup>c</sup>,  
-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>g</sup>, -S(=O)<sub>2</sub>N(R<sup>g</sup>)C(=O)OR<sup>h</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>g</sup>,  
-S(=O)<sub>2</sub>N(R<sup>g</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>g</sup>, -NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>g</sup>)C(=O)R<sup>c</sup>,  
-N(R<sup>a</sup>)C(=O)R<sup>g</sup>, -N(R<sup>g</sup>)C(=O)OR<sup>h</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>g</sup>, -N(R<sup>g</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>,  
-N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>g</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>h</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>g</sup>)S(=O)<sub>2</sub>R<sup>c</sup>,  
-N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>g</sup>, -N(R<sup>g</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>g</sup>, -NR<sup>h</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>g</sup>,  
-NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>g</sup>, -NR<sup>g</sup>C<sub>2-6</sub>alkylOR<sup>h</sup> and -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>g</sup>, wherein R<sup>4</sup> is not  
imidazole or any substituted derivative thereof;

R<sup>5</sup> is H, R<sup>c</sup>, C<sub>1-4</sub>haloalkyl, halo, cyano, -C(=O)R<sup>c</sup>, -C(=O)OR<sup>c</sup>, -C(=O)NR<sup>c</sup>R<sup>a</sup>,  
-C(=NR<sup>a</sup>)NR<sup>h</sup>R<sup>a</sup>, -OR<sup>h</sup>, -OC(=O)R<sup>c</sup>, -OC(=O)NR<sup>a</sup>R<sup>h</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>c</sup>,  
-OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>h</sup>, -OC<sub>2-6</sub>alkylOR<sup>h</sup>, -SR<sup>h</sup>, -S(=O)R<sup>c</sup>, -S(=O)<sub>2</sub>R<sup>c</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>,  
-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>c</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>c</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>R<sup>h</sup>,  
-N(R<sup>a</sup>)C(=O)R<sup>c</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>c</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>h</sup>,  
-N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>c</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>h</sup>,  
-C(=O)R<sup>g</sup>, -C(=O)OR<sup>g</sup>, -C(=O)NR<sup>g</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>g</sup>, -OR<sup>g</sup>, -OC(=O)R<sup>g</sup>,  
-OC(=O)NR<sup>a</sup>R<sup>g</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>g</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>g</sup>, -OC<sub>2-6</sub>alkylOR<sup>g</sup>, -SR<sup>g</sup>,  
-S(=O)R<sup>g</sup>, -S(=O)<sub>2</sub>R<sup>g</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>g</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>g</sup>,  
-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>g</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>g</sup>, -NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>a</sup>)C(=O)R<sup>g</sup>,  
-N(R<sup>a</sup>)C(=O)OR<sup>g</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>g</sup>,  
-N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>g</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>g</sup> or -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>g</sup>; or R<sup>5</sup> is a saturated,  
partially saturated or unsaturated 5-, 6- or 7-membered monocyclic or 6-, 7-, 8-, 9-,  
10- or 11-membered bicyclic ring containing 0, 1, 2, 3 or 4 atoms selected from N, O  
and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups  
and the ring is substituted by 0, 1, 2, 3 or 4 substituents selected from R<sup>c</sup>,  
C<sub>1-4</sub>haloalkyl, halo, cyano, nitro, -C(=O)R<sup>c</sup>, -C(=O)OR<sup>c</sup>, -C(=O)NR<sup>c</sup>R<sup>a</sup>,  
-C(=NR<sup>a</sup>)NR<sup>h</sup>R<sup>a</sup>, -OR<sup>h</sup>, -OC(=O)R<sup>c</sup>, -OC(=O)NR<sup>a</sup>R<sup>h</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>c</sup>,  
-OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>h</sup>, -OC<sub>2-6</sub>alkylOR<sup>h</sup>, -SR<sup>h</sup>, -S(=O)R<sup>c</sup>, -S(=O)<sub>2</sub>R<sup>c</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>,

-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>c</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>c</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>R<sup>h</sup>,  
-N(R<sup>a</sup>)C(=O)R<sup>c</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>c</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>h</sup>,  
-N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>c</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>h</sup>,  
-C(=O)R<sup>g</sup>, -C(=O)OR<sup>g</sup>, -C(=O)NR<sup>g</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>g</sup>, -OR<sup>g</sup>, -OC(=O)R<sup>g</sup>,  
-OC(=O)NR<sup>a</sup>R<sup>g</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>g</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>g</sup>, -OC<sub>2-6</sub>alkylOR<sup>g</sup>, -SR<sup>g</sup>,  
-S(=O)R<sup>g</sup>, -S(=O)<sub>2</sub>R<sup>g</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>g</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>g</sup>,  
-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>g</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>g</sup>, -NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>a</sup>)C(=O)R<sup>g</sup>,  
-N(R<sup>a</sup>)C(=O)OR<sup>g</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>g</sup>,  
-N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>g</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>g</sup> and -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>g</sup>;

R<sup>5'</sup> is H, R<sup>c</sup>, C<sub>1-4</sub>haloalkyl, halo, cyano, -C(=O)R<sup>c</sup>, -C(=O)OR<sup>c</sup>, -C(=O)NR<sup>c</sup>R<sup>a</sup>,  
-C(=NR<sup>a</sup>)NR<sup>h</sup>R<sup>a</sup>, -OR<sup>h</sup>, -OC(=O)R<sup>c</sup>, -OC(=O)NR<sup>a</sup>R<sup>h</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>c</sup>,  
-OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>h</sup>, -OC<sub>2-6</sub>alkylOR<sup>h</sup>, -SR<sup>h</sup>, -S(=O)R<sup>c</sup>, -S(=O)<sub>2</sub>R<sup>c</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>,  
-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>c</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>c</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>R<sup>h</sup>,  
-N(R<sup>a</sup>)C(=O)R<sup>c</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>c</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>h</sup>,  
-N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>c</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>h</sup>,  
-C(=O)R<sup>g</sup>, -C(=O)OR<sup>g</sup>, -C(=O)NR<sup>g</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>g</sup>, -OR<sup>g</sup>, -OC(=O)R<sup>g</sup>,  
-OC(=O)NR<sup>a</sup>R<sup>g</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>g</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>g</sup>, -OC<sub>2-6</sub>alkylOR<sup>g</sup>, -SR<sup>g</sup>,  
-S(=O)R<sup>g</sup>, -S(=O)<sub>2</sub>R<sup>g</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>g</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>g</sup>,  
-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>g</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>g</sup>, -NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>a</sup>)C(=O)R<sup>g</sup>,  
-N(R<sup>a</sup>)C(=O)OR<sup>g</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>g</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>g</sup>,  
-N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>g</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>g</sup> or -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>g</sup>; or R<sup>5'</sup> is a saturated,  
partially saturated or unsaturated 5-, 6- or 7-membered monocyclic or 6-, 7-, 8-, 9-,  
10- or 11-membered bicyclic ring containing 0, 1, 2, 3 or 4 atoms selected from N, O  
and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups  
and the ring is substituted by 0, 1, 2, 3 or 4 substituents selected from R<sup>c</sup>,

C<sub>1-4</sub>haloalkyl, halo, cyano, nitro, -C(=O)R<sup>c</sup>, -C(=O)OR<sup>c</sup>, -C(=O)NR<sup>c</sup>R<sup>a</sup>,  
-C(=NR<sup>a</sup>)NR<sup>h</sup>R<sup>a</sup>, -OR<sup>h</sup>, -OC(=O)R<sup>c</sup>, -OC(=O)NR<sup>a</sup>R<sup>h</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>c</sup>,  
-OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>h</sup>, -OC<sub>2-6</sub>alkylOR<sup>h</sup>, -SR<sup>h</sup>, -S(=O)R<sup>c</sup>, -S(=O)<sub>2</sub>R<sup>c</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>,  
-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>c</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>c</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>R<sup>h</sup>,  
-N(R<sup>a</sup>)C(=O)R<sup>c</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>c</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>h</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>h</sup>,  
-N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>c</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>h</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>h</sup>,  
-C(=O)R<sup>g</sup>, -C(=O)OR<sup>g</sup>, -C(=O)NR<sup>g</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>g</sup>, -OR<sup>g</sup>, -OC(=O)R<sup>g</sup>,  
-OC(=O)NR<sup>a</sup>R<sup>g</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>g</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>g</sup>, -OC<sub>2-6</sub>alkylOR<sup>g</sup>, -SR<sup>g</sup>,

$-S(=O)R^g$ ,  $-S(=O)_2R^g$ ,  $-S(=O)_2NR^aR^g$ ,  $-S(=O)_2N(R^a)C(=O)R^g$ ,  
 $-S(=O)_2N(R^a)C(=O)OR^g$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^g$ ,  $-NR^aR^g$ ,  $-N(R^a)C(=O)R^g$ ,  
 $-N(R^a)C(=O)OR^g$ ,  $-N(R^a)C(=O)NR^aR^g$ ,  $-N(R^a)C(=NR^a)NR^aR^g$ ,  $-N(R^a)S(=O)_2R^g$ ,  
 $-N(R^a)S(=O)_2NR^aR^g$ ,  $-NR^aC_{2-6}alkylNR^aR^g$  and  $-NR^aC_{2-6}alkylOR^g$ ;

$R^6$  is H,  $C_{1-4}haloalkyl$ , halo, cyano,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  
 $-C(=NR^a)NR^aR^a$ ,  $-OH$ ,  $-OC_{2-6}alkyl$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  
 $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  
 $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  
 $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^b$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  
 $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-N(R^a)S(=O)_2NR^aR^a$ ,  
 $-NR^aC_{2-6}alkylNR^aR^a$ ,  $-NR^aC_{2-6}alkylOR^a$  and  $R^c$ ;

$R^{6'}$  is H,  $C_{1-4}haloalkyl$ , halo, cyano,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  
 $-C(=NR^a)NR^aR^a$ ,  $-OH$ ,  $-OC_{2-6}alkyl$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  
 $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  
 $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  
 $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^b$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  
 $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-N(R^a)S(=O)_2NR^aR^a$ ,  
 $-NR^aC_{2-6}alkylNR^aR^a$ ,  $-NR^aC_{2-6}alkylOR^a$  and  $R^c$ ; wherein at least one of  $R^6$  and  $R^{6'}$  is other than H;

$R^a$  is independently, at each instance, H or  $R^b$ ;

$R^b$  is independently, at each instance, phenyl, benzyl or  $C_{1-6}alkyl$ , the phenyl, benzyl and  $C_{1-6}alkyl$  being substituted by 0, 1, 2 or 3 substituents selected from halo,  $C_{1-4}alkyl$ ,  $C_{1-3}haloalkyl$ ,  $-OC_{1-4}alkyl$ ,  $-NH_2$ ,  $-NHC_{1-4}alkyl$ ,  $-N(C_{1-4}alkyl)C_{1-4}alkyl$ ;

$R^c$  is independently at each instance a saturated, partially saturated or unsaturated 5-, 6- or 7-membered monocyclic or 6-, 7-, 8-, 9-, 10- or 11-membered bicyclic ring containing 1, 2, 3 or 4 atoms selected from N, O and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups;

$R^d$  is independently at each instance  $C_{1-8}alkyl$ ,  $C_{1-4}haloalkyl$ , halo, cyano, nitro,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OR^a$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^a$ ,  $-N(R^a)C(=O)R^b$ ,

$-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  
 $-N(R^a)S(=O)_2NR^aR^a$ ,  $-NR^aC_{2-6}alkylNR^aR^a$  or  $-NR^aC_{2-6}alkylOR^a$ ;

$R^c$  is independently at each instance  $C_{1-6}alkyl$  substituted by 0, 1, 2 or 3 substituents independently selected from  $R^d$  and additionally substituted by 0 or 1 substituents selected from  $R^g$ ;

$R^g$  is independently at each instance a saturated, partially saturated or unsaturated 5-, 6- or 7-membered monocyclic or 6-, 7-, 8-, 9-, 10- or 11-membered bicyclic ring containing 0, 1, 2, 3 or 4 atoms selected from N, O and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups and the ring is substituted by 0, 1, 2 or 3 substituents selected from  $C_{1-8}alkyl$ ,  $C_{1-4}haloalkyl$ , halo, cyano, nitro,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OR^a$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^a$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-N(R^a)S(=O)_2NR^aR^a$ ,  $-NR^aC_{2-6}alkylNR^aR^a$  and  $-NR^aC_{2-6}alkylOR^a$ ; and

$R^h$  is independently at each instance  $R^e$  or H.

Claim 2 (canceled)

Claim 3 (currently amended): The compound according to Claim 1, wherein  $Q^3$  is  $C(R^5)$ ,  ~~$Q^4$  is  $C(R^6)$ ,  $Q^5$  is  $C(R^6)$~~ , and  $Q^6$  is  $C(R^5)$ .

Claim 4 (original): The compound according to Claim 3, wherein  $R^4$  is a ring selected from thiophene, pyrrole, 1,3-oxazole, 1,3-thiazole, 1,3,4-oxadiazole, 1,3,4-thiadiazole, 1,2,3-oxadiazole, 1,2,3-thiadiazole, 1H-1,2,3-triazole, isothiazole, 1,2,4-oxadiazole, 1,2,4-thiadiazole, 1,2,3,4-oxatriazole, 1,2,3,4-thiatriazole, 1H-1,2,3,4-tetraazole, 1,2,3,5-oxatriazole, 1,2,3,5-thiatriazole, furan, imidazol-1-yl, imidazol-4-yl, 1,2,4-triazol-4-yl, 1,2,4-triazol-5-yl, isoxazol-3-yl, isoxazol-5-yl, thiolane, pyrrolidine, tetrahydrofuran, 4,5-dihydrothiophene, 2-pyrroline, 4,5-dihydrofuran, pyridazine, pyrimidine, pyrazine, 1,2,3-triazine, 1,2,4-triazine, 1,2,4-triazine, 1,3,5-triazine, pyridine, 2H-3,4,5,6-tetrahydropyran, thiane, 1,2-diazaperhydroine, 1,3-diazaperhydroine, piperazine, 1,3-oxazaperhydroine, morpholine, 1,3-

thiazaperhydroine, 1,4-thiazaperhydroine, piperidine, 2H-3,4-dihydropyran, 2,3-dihydro-4H-thiin, 1,4,5,6-tetrahydropyridine, 2H-5,6-dihydropyran, 2,3-dihydro-6H-thiin, 1,2,5,6-tetrahydropyridine, 3,4,5,6-tetrahydropyridine, 4H-pyran, 4H-thiin, 1,4-dihydropyridine, 1,4-dithiane, 1,4-dioxane, 1,4-oxathiane, 1,2-oxazolidine, 1,2-thiazolidine, pyrazolidine, 1,3-oxazolidine, 1,3-thiazolidine, imidazolidine, 1,2,4-oxadiazolidine, 1,3,4-oxadiazolidine, 1,2,4-thiadiazolidine, 1,3,4-thiadiazolidine, 1,2,4-triazolidine, 2-imidazoline, 3-imidazoline, 2-pyrazoline, 4-imidazoline, 2,3-dihydroisothiazole, 4,5-dihydroisoxazole, 4,5-dihydroisothiazole, 2,5-dihydroisoxazole, 2,5-dihydroisothiazole, 2,3-dihydroisoxazole, 4,5-dihydrooxazole, 2,3-dihydrooxazole, 2,5-dihydrooxazole, 4,5-dihydrothiazole, 2,3-dihydrothiazole,, 2,5-dihydrothiazole, 1,3,4-oxathiazolidine, 1,4,2-oxathiazolidine, 2,3-dihydro-1H-[1,2,3]triazole, 2,5-dihydro-1H-[1,2,3]triazole, 4,5-dihydro-1H-[1,2,3]triazole, 2,3-dihydro-1H-[1,2,4]triazole, 4,5-dihydro-1H-[1,2,4]triazole, 2,3-dihydro-[1,2,4]oxadiazole, 2,5-dihydro-[1,2,4]oxadiazole, 4,5-dihydro-[1,2,4]thiadiazole, 2,3-dihydro-[1,2,4] thidiazole, 2,5-dihydro-[1,2,4] thiadiazole, 4,5-dihydro-[1,2,4] thiadiazole, 2,5-dihydro-[1,2,4]oxadiazole, 2,3-dihydro-[1,2,4]oxadiazole, 4,5-dihydro-[1,2,4]oxadiazole, 2,5-dihydro-[1,2,4]thiadiazole, 2,3-dihydro-[1,2,4] thiadiazole, 4,5-dihydro-[1,2,4] thiadiazole, 2,3-dihydro-[1,3,4]oxadiazole, 2,3-dihydro-[1,3,4]thiadiazole, [1,4,2]oxathiazole, [1,3,4]oxathiazole, 1,3,5-triazaperhydroine, 1,2,4-triazaperhydroine, 1,4,2-dithiazaperhydroine, 1,4,2-dioxaperhydroine, 1,3,5-oxadiazaperhydroine, 1,2,5-oxadiazaperhydroine, 1,3,4-thiadiazaperhydroine, 1,3,5-thiadiazaperhydroine, 1,2,5-thiadiazaperhydroine, 1,3,4-oxadiazaperhydroine, 1,4,3-oxathiazaperhydroine, 1,4,2-oxathiazaperhydroine, 1,4,5,6-tetrahydropyridazine, 1,2,3,4-tetrahydropyridazine, 1,2,3,6-tetrahydropyridazine, 1,2,5,6-tetrahydropyrimidine, 1,2,3,4-tetrahydropyrimidine, 1,4,5,6-tetrahydropyrimidine, 1,2,3,6-tetrahydropyrazine, 1,2,3,4-tetrahydropyrazine, 5,6-dihydro-4H-[1,2]oxazine, 5,6-dihydro-2H-[1,2]oxazine, 3,6-dihydro-2H-[1,2]oxazine, 3,4-dihydro-2H-[1,2]oxazine, 5,6-dihydro-4H-[1,2]thiazine, 5,6-dihydro-2H-[1,2] thiazine, 3,6-dihydro-2H-[1,2] thiazine, 3,4-dihydro-2H-[1,2] thiazine, 5,6-dihydro-2H-[1,3]oxazine, 5,6-dihydro-4H-[1,3]oxazine, 3,6-dihydro-2H-[1,3]oxazine, 3,4-dihydro-2H-[1,3]oxazine, 3,6-dihydro-2H-[1,4]oxazine, 3,4-dihydro-2H-[1,4]oxazine, 5,6-dihydro-2H-[1,3]thiazine, 5,6-dihydro-4H-[1,3]thiazine, 3,6-dihydro-2H-[1,3]thiazine, 3,4-dihydro-2H-[1,3]thiazine, 3,6-dihydro-2H-

[1,4]thiazine, 3,4-dihydro-2H-[1,4]thiazine, 1,2,3,6-tetrahydro-[1,2,4]triazine, 1,2,3,4-tetrahydro-[1,2,4]triazine, 1,2,3,4-tetrahydro-[1,3,5]triazine, 2,3,4,5-tetrahydro-[1,2,4]triazine, 1,4,5,6-tetrahydro-[1,2,4]triazine, 5,6-dihydro-[1,4,2]dioxazine, 5,6-dihydro-[1,4,2]dioxazine, 5,6-dihydro-[1,4,2]dithiazine, 2,3-dihydro-[1,4,2]dioxazine, 3,4-dihydro-2H-[1,3,4]oxadiazine, 3,6-dihydro-2H-[1,3,4]oxadiazine, 3,4-dihydro-2H-[1,3,5]oxadiazine, 3,6-dihydro-2H-[1,3,5]oxadiazine, 5,6-dihydro-2H-[1,2,5]oxadiazine, 5,6-dihydro-4H-[1,2,5]oxadiazine, 3,4-dihydro-2H-[1,3,4]thiadiazine, 3,6-dihydro-2H-[1,3,4]thiadiazine, 3,4-dihydro-2H-[1,3,5]thiadiazine, 3,6-dihydro-2H-[1,3,5]thiadiazine, 5,6-dihydro-2H-[1,2,5]thiadiazine, 5,6-dihydro-4H-[1,2,5]thiadiazine, 5,6-dihydro-2H-[1,2,3]oxadiazine, 3,6-dihydro-2H-[1,2,5]oxadiazine, 5,6-dihydro-4H-[1,3,4]oxadiazine, 3,4-dihydro-2H-[1,2,5]oxadiazine, 5,6-dihydro-2H-[1,2,3]thiadiazine, 3,6-dihydro-2H-[1,2,5]thiadiazine, 5,6-dihydro-4H-[1,3,4]thiadiazine, 3,4-dihydro-2H-[1,2,5]thiadiazine, 5,6-dihydro-[1,4,3]oxathiazine, 5,6-dihydro-[1,4,2]oxathiazine, 2,3-dihydro-[1,4,3]oxathiazine, 2,3-dihydro-[1,4,2]oxathiazine, 4,5-dihydropyridine, 1,6-dihydropyridine, 5,6-dihydropyridine, 2H-pyran, 2H-thiin, 3,6-dihydropyridine, 2,3-dihydropyridazine, 2,5-dihydropyridazine, 4,5-dihydropyridazine, 1,2-dihydropyridazine, 2,3-dihydropyrimidine, 2,5-dihydropyrimidine, 5,6-dihydropyrimidine, 3,6-dihydropyrimidine, 4,5-dihydropyrazine, 5,6-dihydropyrazine, 3,6-dihydropyrazine, 4,5-dihydropyrazine, 1,4-dihydropyrazine, 1,4-dithiin, 1,4-dioxin, 2H-1,2-oxazine, 6H-1,2-oxazine, 4H-1,2-oxazine, 2H-1,3-oxazine, 4H-1,3-oxazine, 6H-1,3-oxazine, 2H-1,4-oxazine, 4H-1,4-oxazine, 2H-1,3-thiazine, 2H-1,4-thiazine, 4H-1,2-thiazine, 6H-1,3-thiazine, 4H-1,4-thiazine, 2H-1,2-thiazine, 6H-1,2-thiazine, 1,4-oxathiin, 2H,5H-1,2,3-triazine, 1H,4H-1,2,3-triazine, 4,5-dihydro-1,2,3-triazine, 1H,6H-1,2,3-triazine, 1,2-dihydro-1,2,3-triazine, 2,3-dihydro-1,2,4-triazine, 3H,6H-1,2,4-triazine, 1H,6H-1,2,4-triazine, 3,4-dihydro-1,2,4-triazine, 1H,4H-1,2,4-triazine, 5,6-dihydro-1,2,4-triazine, 4,5-dihydro-1,2,4-triazine, 2H,5H-1,2,4-triazine, 1,2-dihydro-1,2,4-triazine, 1H,4H-1,3,5-triazine, 1,2-dihydro-1,3,5-triazine, 1,4,2-dithiazine, 1,4,2-dioxazine, 2H-1,3,4-oxadiazine, 2H-1,3,5-oxadiazine, 6H-1,2,5-oxadiazine, 4H-1,3,4-oxadiazine, 4H-1,3,5-oxadiazine, 4H-1,2,5-oxadiazine, 2H-1,3,5-thiadiazine, 6H-1,2,5-thiadiazine, 4H-1,3,4-thiadiazine, 4H-1,3,5-thiadiazine, 4H-1,2,5-thiadiazine, 2H-1,3,4-thiadiazine, 6H-1,3,4-thiadiazine, 6H-1,3,4-oxadiazine, 1,4,2-oxathiazine

and any bicyclic derivative of any of the above rings containing a vicinally-fused phenyl, pyridine or pyrimidine, wherein the carbon atoms of the ring and bicyclic derivative are substituted by 0, 1 or 2 oxo or thioxo groups; wherein the ring or bicyclic derivative there of is substituted by 0, 1, 2, 3 or 4 substituents selected from  $R^c$ ,  $R^e$ , halo,  $C_{1-4}$ haloalkyl, cyano, nitro,  $-C(=O)R^e$ ,  $-C(=O)OR^h$ ,  $-C(=O)NR^aR^h$ ,  $-C(=NR^a)NR^aR^h$ ,  $-OR^h$ ,  $-OC(=O)R^e$ ,  $-OC(=O)NR^aR^h$ ,  $-OC(=O)N(R^a)S(=O)_2R^e$ ,  $-OC_{2-6}alkylNR^aR^h$ ,  $-OC_{2-6}alkylOR^h$ ,  $-SR^e$ ,  $-S(=O)R^e$ ,  $-S(=O)_2R^e$ ,  $-S(=O)_2NR^aR^h$ ,  $-S(=O)_2N(R^a)C(=O)R^e$ ,  $-S(=O)_2N(R^a)C(=O)OR^h$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^h$ ,  $-NR^aR^h$ ,  $-N(R^a)C(=O)R^e$ ,  $-N(R^a)C(=O)OR^h$ ,  $-N(R^a)C(=O)NR^aR^h$ ,  $-N(R^a)C(=NR^a)NR^aR^h$ ,  $-N(R^a)S(=O)_2R^e$ ,  $-N(R^a)S(=O)_2NR^aR^h$ ,  $-NR^aC_{2-6}alkylNR^aR^h$ ,  $-NR^aC_{2-6}alkylOR^h$ ,  $-C(=O)R^g$ ,  $-C(=O)OR^g$ ,  $-C(=O)NR^aR^g$ ,  $-C(=NR^a)NR^aR^g$ ,  $-OR^g$ ,  $-OC(=O)R^g$ ,  $-OC(=O)NR^aR^g$ ,  $-OC(=O)N(R^a)S(=O)_2R^g$ ,  $-OC(=O)N(R^g)S(=O)_2R^e$ ,  $-OC_{2-6}alkylNR^aR^g$ ,  $-OC_{2-6}alkylOR^g$ ,  $-SR^g$ ,  $-S(=O)R^g$ ,  $-S(=O)_2R^g$ ,  $-S(=O)_2NR^aR^g$ ,  $-S(=O)_2N(R^g)C(=O)R^e$ ,  $-S(=O)_2N(R^a)C(=O)R^g$ ,  $-S(=O)_2N(R^g)C(=O)OR^h$ ,  $-S(=O)_2N(R^a)C(=O)OR^g$ ,  $-S(=O)_2N(R^g)C(=O)NR^aR^h$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^g$ ,  $-NR^aR^g$ ,  $-N(R^g)C(=O)R^e$ ,  $-N(R^a)C(=O)R^g$ ,  $-N(R^g)C(=O)OR^h$ ,  $-N(R^a)C(=O)OR^g$ ,  $-N(R^g)C(=O)NR^aR^h$ ,  $-N(R^a)C(=O)NR^aR^g$ ,  $-N(R^g)C(=NR^a)NR^aR^h$ ,  $-N(R^a)C(=NR^a)NR^aR^g$ ,  $-N(R^g)S(=O)_2R^e$ ,  $-N(R^a)S(=O)_2R^g$ ,  $-N(R^g)S(=O)_2NR^aR^h$ ,  $-N(R^a)S(=O)_2NR^aR^g$ ,  $-NR^hC_{2-6}alkylNR^aR^g$ ,  $-NR^aC_{2-6}alkylNR^aR^g$ ,  $-NR^gC_{2-6}alkylOR^h$  and  $-NR^aC_{2-6}alkylOR^g$ .

Claim 5 (original): The compound according to Claim 4, wherein  $R^5$  is  $R^e$ .

Claim 6 (original): The compound according to Claim 4, wherein  $R^{5'}$  is  $R^e$ .

Claim 7 (original): The compound according to Claim 4, wherein  $R^6$  is selected from  $C_{1-4}$ haloalkyl, halo, cyano,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OH$ ,  $-OC_{2-6}alkyl$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^b$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-N(R^a)S(=O)_2NR^aR^a$ ,  $-NR^aC_{2-6}alkylNR^aR^a$ ,  $-NR^aC_{2-6}alkylOR^a$  and  $R^e$ .



Claim 8 (original): The compound according to Claim 7, wherein R<sup>6</sup> is selected from C<sub>1-4</sub>haloalkyl, halo and R<sup>e</sup>.

Claim 9 (original): The compound according to Claim 7, wherein R<sup>6</sup> is selected from C<sub>1-4</sub>haloalkyl and C<sub>2-5</sub>alkyl.

Claim 10 (original): The compound according to Claim 7, wherein R<sup>6</sup> is selected from trifluoromethyl and tert-butyl.

Claim 11 (currently amended): ~~The A compound according to Claim 1, wherein the compound is selected from:~~

7-pyridin-2-yl-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;

7-(1,3-thiazol-2-yl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;

7-pyrazin-2-yl-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;

2-[4-(3,5-dichloropyridin-2-yl)piperazin-1-yl]-7-(3,4-difluorophenyl)-5-(trifluoromethyl)-1H-benzimidazole;

7-bromo-2-[4-(3,5-difluoropyridin-2-yl)piperazin-1-yl]-5-(trifluoromethyl)-1H-benzimidazole;

2-[4-(3,5-dichloropyridin-2-yl)piperazin-1-yl]-5-(trifluoromethyl)-7-[3-(trifluoromethyl)phenyl]-1H-benzimidazole;

2-[4-(3-bromopyridin-2-yl)piperazin-1-yl]-5,7-bis(trifluoromethyl)-1H-benzimidazole;

7-(3,4-difluorobenzyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;

2-[(2R)-4-(3-bromopyridin-2-yl)-2-methylpiperazin-1-yl]-7-chloro-5-(trifluoromethyl)-1H-benzimidazole;

2-[(3R)-4-(3-bromopyridin-2-yl)-3-methylpiperazin-1-yl]-5,7-bis(trifluoromethyl)-1H-benzimidazole;

(2-{4-[7-chloro-5-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)methanol;

(2-{4-[7-bromo-5-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)methanol;

7-chloro-2-{(3R)-3-methyl-4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-5-(trifluoromethyl)-1H-benzimidazole;

7-chloro-2-[(3R)-4-(3-chloropyridin-2-yl)-3-methylpiperazin-1-yl]-5-(trifluoromethyl)-1H-benzimidazole;

7-bromo-2-[(3R)-4-(3-bromopyridin-2-yl)-3-methylpiperazin-1-yl]-5-(trifluoromethyl)-1H-benzimidazole;

2-[(3R)-4-(3-bromopyridin-2-yl)-3-methylpiperazin-1-yl]-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;

2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;

2-[4-(3,5-dichloropyridin-2-yl)piperazin-1-yl]-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;

2-[4-(3-bromopyridin-2-yl)piperazin-1-yl]-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;

N-methyl-2-{4-[5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridine-3-sulfonamide;

7-nitro-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;

5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-amine;

N-(3,4,5-trifluorobenzyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-amine;

(5-chloro-6-{4-[5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)methanol;

(5-chloro-6-{4-[7-iodo-5-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)methanol;

(5-chloro-6-{(3R)-4-[7-iodo-5-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl}pyridin-3-yl)methanol;

7-iodo-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;

4-[(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)amino]butan-2-ol;

3-hydroxy-2,2-dimethyl-N-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)propanamide;

3-hydroxy-N-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)butanamide;

2-[(2R)-4-(3-bromopyridin-2-yl)-2-methylpiperazin-1-yl]-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;

2-[(2S)-4-(3-bromopyridin-2-yl)-2-methylpiperazin-1-yl]-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;

3,5-difluoro-6-{4-[5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-2-amine;

2,2-dimethyl-N-[3-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)pyridin-2-yl]propanamide;

3-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)pyridin-2-amine;

(5-chloro-6-{4-[7-(3-fluoro-4-methoxyphenyl)-5-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)methanol;

2-[(3S)-4-(3-bromopyridin-2-yl)-3-methylpiperazin-1-yl]-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;

2-[(2R)-4-(3-chloropyridin-2-yl)-2-methylpiperazin-1-yl]-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;

4-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)benzylamine;

N-isopropyl-N-[4-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)benzyl]amine;

7-bromo-2-[(2R)-4-(3-chloropyridin-2-yl)-2-methylpiperazin-1-yl]-5-(trifluoromethyl)-1H-benzimidazole;

6-trifluoromethyl-2-[4-(3-trifluoromethyl-pyridin-2-yl)-piperazin-1-yl]-4-(3,4,5-trifluoro-phenyl)-1H-benzoimidazole;

6-trifluoromethyl-2-[4-(3-trifluoromethyl-3,4,5,6-tetrahydro-pyridin-2-yl)-piperazin-1-yl]-4-(3,4,5-trifluoro-phenyl)-1H-benzoimidazole;  
2-[(2R)-4-(3-chloropyridin-2-yl)-2-methylpiperazin-1-yl]-7-(3-fluoro-4-methoxyphenyl)-5-(trifluoromethyl)-1H-benzimidazole;  
7-bromo-2-{(2R)-2-methyl-4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-5-(trifluoromethyl)-1H-benzimidazole;  
2-{(2R)-2-methyl-4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;  
N,N-dimethyl-4-[2-{(2R)-2-methyl-4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-5-(trifluoromethyl)-1H-benzimidazol-7-yl]aniline;  
1-benzyl-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
1-benzyl-6-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
5-chloro-6-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
5-chloro-2-[4-(3,5-dichloropyridin-2-yl)piperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;  
5-chloro-2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;  
6-chloro-5-methyl-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
6-(3,4-difluorophenyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
5-bromo-6-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
5-bromo-2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;  
5-(3,4-difluorophenyl)-6-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
2-[4-(5-chloropyrimidin-4-yl)piperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;  
5-bromo-2-[4-(5-chloropyrimidin-4-yl)piperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;

methyl 5-chloro-6-{4-[6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}nicotinate;  
5-chloro-N-methyl-6-{4-[6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}nicotinamide;  
5-chloro-6-{4-[6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}nicotinamide;  
1-(5-chloro-6-{4-[6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)ethanone;  
1-(5-chloro-6-{4-[6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)ethanol;  
6-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-3H-imidazo[4,5-b]pyridine;  
1-[5-(trifluoromethyl)-1H-benzimidazol-2-yl]-4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-2-one;  
2-{2-(methoxymethyl)-4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-5-(trifluoromethyl)-1H-benzimidazole;  
1-[3-(trifluoromethyl)pyridin-2-yl]-4-[5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]-1,2,3,4-tetrahydroquinoxaline;  
2,2-dimethyl-N-[3-({4-[5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}carbonyl)pyridin-4-yl]propanamide;  
2-[4-(2-piperazin-1-ylpropanoyl)piperazin-1-yl]-6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazole;  
2-[4-(pyridin-2-ylmethyl)piperazin-1-yl]-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;  
3-(trifluoromethyl)-1'-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]-1',2',3',6'-tetrahydro-2,4'-bipyridine;  
2-[(2R)-2-butyl-4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl]-5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazole;  
N~1~,N~1~-dimethyl-N~2~-(5-(trifluoromethyl)-6-{4-[5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)ethane-1,2-diamine;  
N-(5-(trifluoromethyl)-6-{4-[5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)acetamide;

5-fluoro-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole-5-  
carbonitrile;  
methyl 2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole-5-  
carboxylate;  
5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-  
imidazo[4,5-b]pyridine;  
5-(trifluoromethyl)-6-[4-(trifluoromethyl)phenyl]-2-{4-[3-(trifluoromethyl)pyridin-2-  
yl]piperazin-1-yl}-1H-benzimidazole;  
5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-6-(3,4,5-  
trifluorophenyl)-1H-benzimidazole;  
(6-{4-[6-bromo-5-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}-5-  
chloropyridin-3-yl)methanol;  
5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-  
benzimidazole-6-carbonitrile;  
tert-butyl 4-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-  
1H-benzimidazol-6-yl)-3,6-dihydropyridine-1(2H)-carboxylate;  
5-(trifluoromethyl)-6-[4-(trifluoromethyl)cyclohex-1-en-1-yl]-2-{4-[3-  
(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
1-{4-[5-(trifluoromethyl)-7-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-  
1-yl}isoquinoline;  
(6-{(3R)-4-[6-bromo-5-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-  
1-yl}-5-chloropyridin-3-yl)methanol;  
6-{4-[7-bromo-5-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}quinoline;  
6-(trifluoromethyl)-N-[4-(trifluoromethyl)benzyl]-2-{4-[3-(trifluoromethyl)pyridin-2-  
yl]piperazin-1-yl}-1H-benzimidazol-5-amine;  
4-{4-[7-bromo-5-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}quinoline;  
N-(2-piperidin-1-ylethyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-  
yl]piperazin-1-yl}-1H-benzimidazol-6-amine;  
N-(2-morpholin-4-ylethyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-  
yl]piperazin-1-yl}-1H-benzimidazol-6-amine;  
7-{4-[7-bromo-5-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}quinoline;

7-{4-[7-bromo-5-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}isoquinoline;  
N-(2-piperazin-1-ylethyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-6-amine;  
5-{4-[4-bromo-6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}isoquinoline;  
6-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole-4-carbonitrile;  
tert-butyl 4-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)-3,6-dihydropyridine-1(2H)-carboxylate;  
7-(4-tert-butylcyclohex-1-en-1-yl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
5-(trifluoromethyl)-7-[4-(trifluoromethyl)cyclohex-1-en-1-yl]-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-(1,2,3,6-tetrahydropyridin-4-yl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-(1-isopropyl-1,2,3,6-tetrahydropyridin-4-yl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
5-(trifluoromethyl)-7-[4-(trifluoromethyl)cyclohexyl]-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
2-{4-[5-bromo-3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-6-(trifluoromethyl)-1H-benzimidazole;  
4-bromo-2-{4-[5-bromo-3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-6-(trifluoromethyl)-1H-benzimidazole;  
N-[6-{4-[4-bromo-6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}-5-(trifluoromethyl)pyridin-2-yl]acetamide;  
7-piperidin-1-yl-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-morpholin-4-yl-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-piperazin-1-yl-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;

4-bromo-2-[(2R)-4-(3,5-dichloropyridin-2-yl)-2-methylpiperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;  
(5-chloro-6-[(3R)-3-methyl-4-[6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl]pyridin-3-yl)methanol;  
(6-[(3R)-4-[4-bromo-6-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl]-5-chloropyridin-3-yl)methanol;  
2-[(2R)-2-methyl-4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;  
2-[(2R)-4-[3-chloro-5-(methoxymethyl)pyridin-2-yl]-2-methylpiperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;  
1-(5-chloro-6-[(3R)-3-methyl-4-[6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl]pyridin-3-yl)ethanol;  
N-[(5-chloro-6-[(3R)-3-methyl-4-[6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl]pyridin-3-yl)methyl]acetamide;  
N-[(5-chloro-6-[(3R)-3-methyl-4-[6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl]pyridin-3-yl)methyl]-N-methylacetamide;  
2-(5-chloro-6-[(3R)-3-methyl-4-[6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl]pyridin-3-yl)propan-2-ol;  
2-[(2R)-4-(3,5-dichloropyridin-2-yl)-2-methylpiperazin-1-yl]-6-(trifluoromethyl)-4-[4-(trifluoromethyl)phenyl]-1H-benzimidazole;  
4-(4-chlorophenyl)-2-[(2R)-4-(3,5-dichloropyridin-2-yl)-2-methylpiperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;  
2-[(2R)-4-(3,5-dichloropyridin-2-yl)-2-methylpiperazin-1-yl]-6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazole;  
[5-chloro-6-[(3R)-3-methyl-4-[6-(trifluoromethyl)-4-[4-(trifluoromethyl)phenyl]-1H-benzimidazol-2-yl]piperazin-1-yl]pyridin-3-yl)methanol;  
(5-chloro-6-[(3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl]pyridin-3-yl)methanol;  
(5-chloro-6-[(3R)-3-methyl-4-[4-pyridin-3-yl]-6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl]pyridin-3-yl)methanol;  
(5-chloro-6-[(3R)-4-[4-(4-chlorophenyl)-6-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl]pyridin-3-yl)methanol;



2-[(2R)-4-(3,5-dichloropyridin-2-yl)-2-methylpiperazin-1-yl]-4-pyridin-3-yl-6-(trifluoromethyl)-1H-benzimidazole;  
1-(5-chloro-6-{(3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)ethanol;  
1-[5-chloro-6-{(3R)-3-methyl-4-{6-(trifluoromethyl)-4-[4-(trifluoromethyl)phenyl]-1H-benzimidazol-2-yl}piperazin-1-yl}pyridin-3-yl]ethanol;  
1-(5-chloro-6-{(3R)-4-[4-(4-chlorophenyl)-6-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl}pyridin-3-yl)ethanol;  
1-(5-chloro-6-{(3R)-3-methyl-4-[4-pyridin-3-yl-6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)ethanol;  
N-[2-{(2R)-4-[3-chloro-5-(hydroxymethyl)pyridin-2-yl]-2-methylpiperazin-1-yl}-5-(trifluoromethyl)-1H-benzimidazol-7-yl]-3,4,5-trifluorobenzamide;  
(5-chloro-6-{(3R)-3-methyl-4-[4-[(3,4,5-trifluorobenzyl)amino]-6-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)methanol;  
4-[3-chloro-5-(hydroxymethyl)pyridin-2-yl]-N-methyl-1-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazine-2-carboxamide;  
(1S)-1-(5-chloro-6-{(3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)ethanol;  
(1R)-1-(5-chloro-6-{(3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)ethanol;  
N-[2-{(2R)-4-[3-chloro-5-(hydroxymethyl)pyridin-2-yl]-2-methylpiperazin-1-yl}-6-(trifluoromethyl)-1H-benzimidazol-4-yl]-4-(trifluoromethyl)benzamide;  
N-[2-{(2R)-4-[3-chloro-5-(hydroxymethyl)pyridin-2-yl]-2-methylpiperazin-1-yl}-6-(trifluoromethyl)-1H-benzimidazol-4-yl]cyclohexanecarboxamide;  
4-chloro-N-[2-{(2R)-4-[3-chloro-5-(hydroxymethyl)pyridin-2-yl]-2-methylpiperazin-1-yl}-6-(trifluoromethyl)-1H-benzimidazol-4-yl]benzamide;  
methyl 5-chloro-6-{(3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}nicotinate;  
(6-{(3R)-4-[7-[bis(cyclohexylmethyl)amino]-5-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl}-5-chloropyridin-3-yl)methanol;  
(5-chloro-6-{(3R)-4-[7-[(cyclohexylmethyl)amino]-5-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl}pyridin-3-yl)methanol;

1-(5-chloro-6-((3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)propan-1-ol;  
tert-butyl 2-((2R)-4-[3-chloro-5-(hydroxymethyl)pyridin-2-yl]-2-methylpiperazin-1-yl)-5-(trifluoromethyl)-1H-benzimidazol-7-ylcarbamate;  
[6-((3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl)-5-(trifluoromethyl)pyridin-3-yl]methanol;  
1-[6-((3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl)-5-(trifluoromethyl)pyridin-3-yl]propan-1-ol;  
1-[6-((3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl)-5-(trifluoromethyl)pyridin-3-yl]prop-2-en-1-ol;  
2-methyl-1-[6-((3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl)-5-(trifluoromethyl)pyridin-3-yl]propan-1-ol;  
1-[6-((3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl)-5-(trifluoromethyl)pyridin-3-yl]ethanol;  
[6-((3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl)-5-(trifluoromethyl)pyridin-3-yl](phenyl)methanol;  
[6-((3R)-3-methyl-4-((6-(trifluoromethyl)-4-[4-(trifluoromethyl)phenyl]-1H-benzimidazol-2-yl}piperazin-1-yl)-5-(trifluoromethyl)pyridin-3-yl]methanol;  
1-[6-((3R)-3-methyl-4-((6-(trifluoromethyl)-4-[4-(trifluoromethyl)phenyl]-1H-benzimidazol-2-yl}piperazin-1-yl)-5-(trifluoromethyl)pyridin-3-yl]propan-1-ol;  
1-[6-((3R)-3-methyl-4-((6-(trifluoromethyl)-4-[4-(trifluoromethyl)phenyl]-1H-benzimidazol-2-yl}piperazin-1-yl)-5-(trifluoromethyl)pyridin-3-yl]prop-2-en-1-ol;  
(5-(trifluoromethyl)-6-((4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}pyridin-3-yl)methanol;  
[5-(trifluoromethyl)-6-((4-[6-(trifluoromethyl)-4-[4-(trifluoromethyl)phenyl]-1H-benzimidazol-2-yl]piperazin-1-yl)pyridin-3-yl]methanol;  
1-[6-((3R)-3-methyl-4-((6-(trifluoromethyl)-4-[4-(trifluoromethyl)phenyl]-1H-benzimidazol-2-yl}piperazin-1-yl)-5-(trifluoromethyl)pyridin-3-yl]ethanol;  
5-chloro-6-((3R)-3-methyl-4-[6-(trifluoromethyl)-4-(3,4,5-trifluorophenyl)-1H-benzimidazol-2-yl]piperazin-1-yl}nicotinic acid;  
2-((2R)-4-(3-chloropyridin-2-yl)-2-methylpiperazin-1-yl)-7-(3,5-difluorophenyl)-5-(trifluoromethyl)-1H-benzimidazole;

7-(3,5-difluorophenyl)-2-[(2R)-4-[3-(3,5-difluorophenyl)pyridin-2-yl]-2-methylpiperazin-1-yl]-5-(trifluoromethyl)-1H-benzimidazole;  
2-[(2R)-4-(3-chloropyridin-2-yl)-2-methylpiperazin-1-yl]-7-(3-fluorophenyl)-5-(trifluoromethyl)-1H-benzimidazole;  
7-(3-fluorophenyl)-2-[(2R)-4-[3-(3-fluorophenyl)pyridin-2-yl]-2-methylpiperazin-1-yl]-5-(trifluoromethyl)-1H-benzimidazole;  
4-[2-[(2R)-4-(3-chloropyridin-2-yl)-2-methylpiperazin-1-yl]-5-(trifluoromethyl)-1H-benzimidazol-7-yl]-N,N-dimethylaniline;  
2-[(2R)-4-(3-chloropyridin-2-yl)-2-methylpiperazin-1-yl]-5-(trifluoromethyl)-7-[4-(trifluoromethyl)phenyl]-1H-benzimidazole;  
2-[(2R)-4-(3-chloropyridin-2-yl)-2-methylpiperazin-1-yl]-7-(4-fluorophenyl)-5-(trifluoromethyl)-1H-benzimidazole;  
5-bromo-2-[4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl]-1H-imidazo[4,5-b]pyridine;  
2-[4-(4-Chloro-[1,2,5]thiadiazol-3-yl)-piperazin-1-yl]-5-trifluoromethyl-7-(3,4,5-trifluoro-phenyl)-1H-benzimidazole;  
6-(trifluoromethyl)-2-[4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl]-1H-benzimidazole;  
N-isopropyl-2-[4-[5-(trifluoromethyl)-1H-benzimidazol-2-yl]piperazin-1-yl]pyridin-3-amine;  
2-[4-(2,4-difluorophenyl)piperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;  
2-[4-(6-chloropyridin-2-yl)piperazin-1-yl]-6-(trifluoromethyl)-1H-benzimidazole;  
7-pyridin-4-yl-5-(trifluoromethyl)-2-[4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl]-1H-benzimidazole;  
7-(3-fluorophenyl)-5-(trifluoromethyl)-2-[4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl]-1H-benzimidazole;  
5-(trifluoromethyl)-7-[2-(trifluoromethyl)phenyl]-2-[4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl]-1H-benzimidazole;  
7-(4-fluorophenyl)-5-(trifluoromethyl)-2-[4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl]-1H-benzimidazole;  
7-(3-chloro-4-fluorophenyl)-5-(trifluoromethyl)-2-[4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl]-1H-benzimidazole;

7-(2-methoxyphenyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-(3,5-difluorophenyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
[3-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)phenyl]methanol;  
7-(6-methoxypyridin-3-yl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
8-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)quinoline;  
7-(4-tert-butylphenyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-(3-chloropyridin-4-yl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-[3-(trifluoromethoxy)phenyl]-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
N-[4-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)phenyl]acetamide;  
7-(1-benzothien-2-yl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
3-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)aniline;  
N,N-dimethyl-4-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)aniline;  
7-(3-fluoro-4-methoxyphenyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-(2,4-dimethoxypyrimidin-5-yl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-(2,5-difluorophenyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-(2,4-difluorophenyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;

7-(3-furyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-(2,3-difluorophenyl)-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
4-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)aniline;  
[4-(5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazol-7-yl)phenyl]methanol;  
7-[(E)-2-phenylvinyl]-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
7-cyclohex-1-en-1-yl-5-(trifluoromethyl)-2-{4-[3-(trifluoromethyl)pyridin-2-yl]piperazin-1-yl}-1H-benzimidazole;  
{5-chloro-6-[(3R)-3-methyl-4-(6-(trifluoromethyl)-4-{[4-(trifluoromethyl)benzyl]amino}-1H-benzimidazol-2-yl)piperazin-1-yl]pyridin-3-yl}methanol;  
(6-{(3R)-4-[7-[bis(4-chlorobenzyl)amino]-5-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl}-5-chloropyridin-3-yl)methanol;  
(5-chloro-6-{(3R)-4-[7-[(4-chlorobenzyl)amino]-5-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl}pyridin-3-yl)methanol;  
(6-{(3R)-4-[7-[bis(4-tert-butylbenzyl)amino]-5-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl}-5-chloropyridin-3-yl)methanol;  
(6-{(3R)-4-[7-[(4-tert-butylbenzyl)amino]-5-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl}-5-chloropyridin-3-yl)methanol;  
(5-chloro-6-{(3R)-4-[7-(dibenzylamino)-5-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl}pyridin-3-yl)methanol; and  
(6-{(3R)-4-[7-(benzylamino)-5-(trifluoromethyl)-1H-benzimidazol-2-yl]-3-methylpiperazin-1-yl}-5-chloropyridin-3-yl)methanol; or a pharmaceutically-acceptable salt thereof.

Claim 12 (canceled)

Claim 13 (previously amended): A pharmaceutical composition comprising a compound according to any one of Claims 1, 3, 4 and 11 and a pharmaceutically-acceptable diluent or carrier.

Claim 14 (canceled)

Claim 15 (new): The compound according to Claim 4, wherein  $R^4$  is a ring selected from pyrimidine and pyridine; wherein the ring is substituted by 0, 1, 2, 3 or 4 substituents selected from  $R^c$ ,  $R^e$ , halo,  $C_{1-4}$ haloalkyl, cyano, nitro,  $-C(=O)R^c$ ,  $-C(=O)OR^h$ ,  $-C(=O)NR^aR^h$ ,  $-C(=NR^a)NR^aR^h$ ,  $-OR^h$ ,  $-OC(=O)R^c$ ,  $-OC(=O)NR^aR^h$ ,  $-OC(=O)N(R^a)S(=O)_2R^c$ ,  $-OC_{2-6}alkylNR^aR^h$ ,  $-OC_{2-6}alkylOR^h$ ,  $-SR^c$ ,  $-S(=O)R^c$ ,  $-S(=O)_2R^c$ ,  $-S(=O)_2NR^aR^h$ ,  $-S(=O)_2N(R^a)C(=O)R^c$ ,  $-S(=O)_2N(R^a)C(=O)OR^h$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^h$ ,  $-NR^aR^h$ ,  $-N(R^a)C(=O)R^c$ ,  $-N(R^a)C(=O)OR^h$ ,  $-N(R^a)C(=O)NR^aR^h$ ,  $-N(R^a)C(=NR^a)NR^aR^h$ ,  $-N(R^a)S(=O)_2R^c$ ,  $-N(R^a)S(=O)_2NR^aR^h$ ,  $-NR^aC_{2-6}alkylNR^aR^h$ ,  $-NR^aC_{2-6}alkylOR^h$ ,  $-C(=O)R^g$ ,  $-C(=O)OR^g$ ,  $-C(=O)NR^aR^g$ ,  $-C(=NR^a)NR^aR^g$ ,  $-OR^g$ ,  $-OC(=O)R^g$ ,  $-OC(=O)NR^aR^g$ ,  $-OC(=O)N(R^a)S(=O)_2R^g$ ,  $-OC(=O)N(R^g)S(=O)_2R^c$ ,  $-OC_{2-6}alkylNR^aR^g$ ,  $-OC_{2-6}alkylOR^g$ ,  $-SR^g$ ,  $-S(=O)R^g$ ,  $-S(=O)_2R^g$ ,  $-S(=O)_2NR^aR^g$ ,  $-S(=O)_2N(R^g)C(=O)R^c$ ,  $-S(=O)_2N(R^a)C(=O)R^g$ ,  $-S(=O)_2N(R^g)C(=O)OR^h$ ,  $-S(=O)_2N(R^a)C(=O)OR^g$ ,  $-S(=O)_2N(R^g)C(=O)NR^aR^h$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^g$ ,  $-NR^aR^g$ ,  $-N(R^g)C(=O)R^c$ ,  $-N(R^a)C(=O)R^g$ ,  $-N(R^g)C(=O)OR^h$ ,  $-N(R^a)C(=O)OR^g$ ,  $-N(R^g)C(=O)NR^aR^h$ ,  $-N(R^a)C(=O)NR^aR^g$ ,  $-N(R^g)C(=NR^a)NR^aR^h$ ,  $-N(R^a)C(=NR^a)NR^aR^g$ ,  $-N(R^g)S(=O)_2R^c$ ,  $-N(R^a)S(=O)_2R^g$ ,  $-N(R^g)S(=O)_2NR^aR^h$ ,  $-N(R^a)S(=O)_2NR^aR^g$ ,  $-NR^hC_{2-6}alkylNR^aR^g$ ,  $-NR^aC_{2-6}alkylNR^aR^g$ ,  $-NR^gC_{2-6}alkylOR^h$  and  $-NR^aC_{2-6}alkylOR^g$ .

Claim 16 (new): The compound according to Claim 4, wherein  $R^4$  is a ring selected from pyrimidine and pyridine; wherein the ring is substituted by 0, 1, 2, 3 or 4 substituents selected from halo and  $C_{1-4}$ haloalkyl.

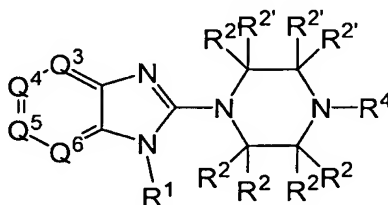
Claim 17 (new): The compound according to Claim 4, wherein  $R^4$  is pyrimidin-4-yl substituted by 1, 2, 3 or 4 substituents selected from halo and  $C_{1-4}$ haloalkyl.

Claim 18 (new): The compound according to Claim 4, wherein R<sup>4</sup> is pyridine substituted by 1, 2, 3 or 4 substituents selected from halo and C<sub>1-4</sub>haloalkyl.

Claim 19 (new): The compound according to Claim 4, wherein R<sup>4</sup> is pyridin-2-yl substituted by 1 or 2 substituents selected from CH<sub>3</sub>, Cl and CF<sub>3</sub>.

Claim 20 (new): The compound according to Claim 4, wherein R<sup>4</sup> is 3-trifluoromethylpyridin-2-yl.

Claim 21 (new): A compound having the structure:



or any pharmaceutically-acceptable salt thereof, wherein:

Q<sup>3</sup> is N or C(R<sup>5</sup>);

Q<sup>4</sup> is N or C(R<sup>6</sup>);

Q<sup>5</sup> is N or C(R<sup>6</sup>);

Q<sup>6</sup> is N or C(R<sup>5</sup>);

R<sup>1</sup> is H or -(C(R<sup>2</sup>)(R<sup>2</sup>))<sub>m</sub>-R<sup>g</sup>;

R<sup>2</sup> is, independently, in each instance, H, C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, -O(C<sub>1-7</sub>alkyl), -N(C<sub>1-7</sub>alkyl)R<sup>a</sup>, or a C<sub>1-6</sub>alkyl substituted by 1, 2 or 3 substituents selected from halo, cyano, -OR<sup>a</sup>, -OC(=O)R<sup>b</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup> and -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup>; wherein any two geminal R<sup>2</sup> groups may additionally be oxo;

R<sup>2'</sup> is, independently, in each instance, H, C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, -O(C<sub>1-7</sub>alkyl), -N(C<sub>1-7</sub>alkyl)R<sup>a</sup>, or a C<sub>1-6</sub>alkyl substituted by 1, 2 or 3 substituents selected from halo, cyano, -OR<sup>a</sup>, -OC(=O)R<sup>b</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup> and -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup>; wherein any two geminal R<sup>2'</sup> groups may additionally be oxo;

$R^4$  is  $R^c$  substituted by 0, 1, 2, 3 or 4 substituents selected from  $C_{1-8}$ alkyl,  $C_{1-4}$ haloalkyl, halo, cyano, nitro,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OR^a$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^a$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-N(R^a)S(=O)_2NR^aR^a$ ,  $-NR^aC_{2-6}alkylNR^aR^a$ ,  $-NR^aC_{2-6}alkylOR^a$  and  $R^e$ ;

$R^5$  is independently, at each instance, H,  $C_{1-8}$ alkyl,  $C_{1-4}$ haloalkyl, halo, cyano,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^bR^a$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OR^a$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^a$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-N(R^a)S(=O)_2NR^aR^a$ ,  $-NR^aC_{2-6}alkylNR^aR^a$ ,  $-NR^aC_{2-6}alkylOR^a$ ,  $C_{1-3}alkylR^c$ ,  $C_{1-3}alkylR^f$  and  $R^e$ ; or  $R^5$  is a saturated or unsaturated 5-, 6- or 7-membered monocyclic or 6-, 7-, 8-, 9-, 10- or 11-membered bicyclic ring containing 0, 1, 2, 3 or 4 atoms selected from N, O and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups and the ring is substituted by 0, 1, 2, 3 or 4 substituents selected from  $C_{1-8}$ alkyl,  $C_{1-4}$ haloalkyl, halo, cyano, nitro,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OR^a$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^a$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-N(R^a)S(=O)_2NR^aR^a$ ,  $-NR^aC_{2-6}alkylNR^aR^a$ ,  $-NR^aC_{2-6}alkylOR^a$  and  $R^e$ ;

$R^6$  is independently, at each instance, H,  $C_{1-8}$ alkyl,  $C_{1-4}$ haloalkyl, halo, cyano,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OH$ ,  $-OC_{2-6}alkyl$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^b$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-N(R^a)S(=O)_2NR^aR^a$ ,  $-NR^aC_{2-6}alkylNR^aR^a$ ,  $-NR^aC_{2-6}alkylOR^a$  and  $R^e$ ;



R<sup>a</sup> is independently, at each instance, H or R<sup>b</sup>;

R<sup>b</sup> is independently, at each instance, phenyl, benzyl or C<sub>1-6</sub>alkyl, the phenyl, benzyl and C<sub>1-6</sub>alkyl being substituted by 0, 1, 2 or 3 substituents selected from halo, C<sub>1-4</sub>alkyl, C<sub>1-3</sub>haloalkyl, -OC<sub>1-4</sub>alkyl, -NH<sub>2</sub>, -NHC<sub>1-4</sub>alkyl, -N(C<sub>1-4</sub>alkyl)C<sub>1-4</sub>alkyl;

R<sup>c</sup> is independently at each instance a saturated or unsaturated 5-, 6- or 7-membered monocyclic or 6-, 7-, 8-, 9-, 10- or 11-membered bicyclic ring containing 1, 2, 3 or 4 atoms selected from N, O and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups;

R<sup>d</sup> is independently at each instance C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, halo, cyano, nitro, -C(=O)R<sup>b</sup>, -C(=O)OR<sup>b</sup>, -C(=O)NR<sup>a</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -OR<sup>a</sup>, -OC(=O)R<sup>b</sup>, -OC(=O)NR<sup>a</sup>R<sup>a</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -OC<sub>2-6</sub>alkylOR<sup>a</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup> or -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup>;

R<sup>e</sup> is independently at each instance C<sub>1-6</sub>alkyl substituted by 1, 2 or 3 substituents independently selected from R<sup>d</sup>;

R<sup>f</sup> is independently at each instance R<sup>c</sup> substituted by 1, 2 or 3 substituents independently selected from R<sup>d</sup>; and

R<sup>g</sup> is independently at each instance a saturated or unsaturated 5-, 6- or 7-membered monocyclic or 6-, 7-, 8-, 9-, 10- or 11-membered bicyclic ring containing 0, 1, 2, 3 or 4 atoms selected from N, O and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups and the ring is substituted by 0, 1, 2 or 3 substituents selected from C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, halo, cyano, nitro, -C(=O)R<sup>b</sup>, -C(=O)OR<sup>b</sup>, -C(=O)NR<sup>a</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -OR<sup>a</sup>, -OC(=O)R<sup>b</sup>, -OC(=O)NR<sup>a</sup>R<sup>a</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -OC<sub>2-6</sub>alkylOR<sup>a</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup> and -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup>.

Claim 22 (new): The compound according to Claim 21, wherein:

Q<sup>3</sup> is C(R<sup>5</sup>);

$Q^4$  is  $C(R^6)$ ;  
 $Q^5$  is  $C(R^6)$ ; and  
 $Q^6$  is  $C(R^5)$ .

Claim 23 (new): The compound according to Claim 21, wherein any one of  $Q^3$ ,  $Q^4$ ,  $Q^5$  and  $Q^6$  is N.

Claim 24 (new): The compound according to Claim 21, wherein  $R^1$  is H.

Claim 25 (new): The compound according to Claim 21, wherein  $R^1$  is  $-(C(R^2)(R^2))_m-R^g$ .

Claim 26 (new): The compound according to Claim 21, wherein  $R^2$  is, in each instance, H.

Claim 27 (new): The compound according to Claim 21, wherein at least one  $R^2$  group is selected from  $C_{1-8}$ alkyl,  $C_{1-4}$ haloalkyl,  $-O(C_{1-7}$ alkyl),  $-N(C_{1-7}$ alkyl) $R^a$ , oxo and  $C_{1-6}$ alkyl substituted by 0, 1, 2 or 3 substituents selected from halo, cyano,  $-OR^a$ ,  $-OC(=O)R^b$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-NR^aR^a$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-NR^aC_{2-6}$ alkyl $NR^aR^a$  and  $-NR^aC_{2-6}$ alkyl $OR^a$ .

Claim 28 (new): The compound according to Claim 21, wherein  $R^4$  is an unsaturated 6- or 7-membered monocyclic or 6-, 7-, 8-, 9-, 10- or 11-membered bicyclic ring containing 1, 2, 3 or 4 atoms selected from N, O and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups and the rings are substituted by 0, 1, 2, 3 or 4 substituents selected from  $C_{1-8}$ alkyl,  $C_{1-4}$ haloalkyl, halo, cyano, nitro,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OR^a$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}$ alkyl $NR^aR^a$ ,  $-OC_{2-6}$ alkyl $OR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^a$ ,  $-N(R^a)C(=O)R^b$ ,

$-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  
 $-N(R^a)S(=O)_2NR^aR^a$ ,  $-NR^aC_{2-6}alkylNR^aR^a$ ,  $-NR^aC_{2-6}alkylOR^a$  and  $R^c$ .

Claim 29 (new): The compound according to Claim 21, wherein  $R^4$  is an unsaturated 6-membered monocyclic ring containing 1, 2 or 3 atoms selected from N, O and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups and the rings are substituted by 0, 1, 2, 3 or 4 substituents selected from  $C_{1-8}alkyl$ ,  $C_{1-4}haloalkyl$ , halo, cyano, nitro,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OR^a$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^a$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-N(R^a)S(=O)_2NR^aR^a$ ,  $-NR^aC_{2-6}alkylNR^aR^a$ ,  $-NR^aC_{2-6}alkylOR^a$  and  $R^c$ .

Claim 30 (new): The compound according to Claim 21, wherein  $R^5$  is, at each instance, H.

Claim 31 (new): The compound according to Claim 21, wherein at least one  $R^5$  is  $C_{1-8}alkyl$ ,  $C_{1-4}haloalkyl$ , halo, cyano,  $-C(=O)R^b$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OR^a$ ,  $-OC(=O)R^b$ ,  $-OC(=O)NR^aR^a$ ,  $-OC(=O)N(R^a)S(=O)_2R^b$ ,  $-OC_{2-6}alkylNR^aR^a$ ,  $-OC_{2-6}alkylOR^a$ ,  $-SR^a$ ,  $-S(=O)R^b$ ,  $-S(=O)_2R^b$ ,  $-S(=O)_2NR^aR^a$ ,  $-S(=O)_2N(R^a)C(=O)R^b$ ,  $-S(=O)_2N(R^a)C(=O)OR^b$ ,  $-S(=O)_2N(R^a)C(=O)NR^aR^a$ ,  $-NR^aR^a$ ,  $-N(R^a)C(=O)R^b$ ,  $-N(R^a)C(=O)OR^b$ ,  $-N(R^a)C(=O)NR^aR^a$ ,  $-N(R^a)C(=NR^a)NR^aR^a$ ,  $-N(R^a)S(=O)_2R^b$ ,  $-N(R^a)S(=O)_2NR^aR^a$ ,  $-NR^aC_{2-6}alkylNR^aR^a$ ,  $-NR^aC_{2-6}alkylOR^a$ ,  $C_{1-3}alkylR^c$ ,  $C_{1-3}alkylR^f$  and  $R^c$ .

Claim 32 (new): The compound according to Claim 21, wherein at least one  $R^5$  is a saturated or unsaturated 5-, 6- or 7-membered monocyclic or 6-, 7-, 8-, 9-, 10- or 11-membered bicyclic ring containing 0, 1, 2, 3 or 4 atoms selected from N, O and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups and the ring is substituted by 0, 1, 2, 3 or 4 substituents selected from  $C_{1-8}alkyl$ ,  $C_{1-4}haloalkyl$ , halo, cyano, nitro,  $-C(=O)R^b$ ,  $-C(=O)OR^b$ ,  $-C(=O)NR^aR^a$ ,  $-C(=NR^a)NR^aR^a$ ,  $-OR^a$ ,

-OC(=O)R<sup>b</sup>, -OC(=O)NR<sup>a</sup>R<sup>a</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>,  
-OC<sub>2-6</sub>alkylOR<sup>a</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>b</sup>,  
-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>,  
-N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>,  
-N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup> and R<sup>c</sup>.

Claim 33 (new): The compound according to Claim 21, wherein at least one R<sup>5</sup> is an unsaturated 5-, 6- or 7-membered monocyclic ring containing 0, 1, 2 or 3 atoms selected from N, O and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups and the ring is substituted by 0, 1, 2, 3 or 4 substituents selected from C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, halo, cyano, nitro, -C(=O)R<sup>b</sup>, -C(=O)OR<sup>b</sup>, -C(=O)NR<sup>a</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -OR<sup>a</sup>, -OC(=O)R<sup>b</sup>, -OC(=O)NR<sup>a</sup>R<sup>a</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -OC<sub>2-6</sub>alkylOR<sup>a</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup> and R<sup>c</sup>.

Claim 34 (new): The compound according to Claim 21, wherein at least one R<sup>5</sup> is an unsaturated 6-membered monocyclic ring containing 0, 1 or 2 N atoms, wherein the ring is substituted by 0, 1, 2, 3 or 4 substituents selected from C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, halo, cyano, nitro, -C(=O)R<sup>b</sup>, -C(=O)OR<sup>b</sup>, -C(=O)NR<sup>a</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -OR<sup>a</sup>, -OC(=O)R<sup>b</sup>, -OC(=O)NR<sup>a</sup>R<sup>a</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -OC<sub>2-6</sub>alkylOR<sup>a</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup> and R<sup>c</sup>.

Claim 35 (new): The compound according to Claim 21, wherein at least one R<sup>5</sup> is a saturated or unsaturated 6-, 7-, 8-, 9-, 10- or 11-membered bicyclic ring containing 0, 1, 2, 3 or 4 atoms selected from N, O and S, wherein the carbon atoms of the ring are substituted by 0, 1 or 2 oxo groups and the ring is substituted by 0, 1, 2, 3 or 4 substituents selected from C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, halo, cyano, nitro, -C(=O)R<sup>b</sup>,

-C(=O)OR<sup>b</sup>, -C(=O)NR<sup>a</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -OR<sup>a</sup>, -OC(=O)R<sup>b</sup>, -OC(=O)NR<sup>a</sup>R<sup>a</sup>,  
-OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -OC<sub>2-6</sub>alkylOR<sup>a</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>,  
-S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>b</sup>,  
-S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>,  
-N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>,  
-NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup> and R<sup>c</sup>.

Claim 36 (new): The compound according to Claim 21, wherein at least one R<sup>6</sup> is selected from C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, halo, cyano, -C(=O)R<sup>b</sup>, -C(=O)OR<sup>b</sup>, -C(=O)NR<sup>a</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -OC(=O)R<sup>b</sup>, -OC(=O)NR<sup>a</sup>R<sup>a</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -OC<sub>2-6</sub>alkylOR<sup>a</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup> and R<sup>c</sup>.

Claim 37 (new): The compound according to Claim 21, wherein R<sup>4</sup> is pyridine or pyrimidine, wherein the pyridine or pyrimidine by substituted by 1, 2, 3 or 4 substituents selected from C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl, halo, cyano, nitro, -C(=O)R<sup>b</sup>, -C(=O)OR<sup>b</sup>, -C(=O)NR<sup>a</sup>R<sup>a</sup>, -C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -OR<sup>a</sup>, -OC(=O)R<sup>b</sup>, -OC(=O)NR<sup>a</sup>R<sup>a</sup>, -OC(=O)N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -OC<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -OC<sub>2-6</sub>alkylOR<sup>a</sup>, -SR<sup>a</sup>, -S(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>R<sup>b</sup>, -S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -S(=O)<sub>2</sub>N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=O)R<sup>b</sup>, -N(R<sup>a</sup>)C(=O)OR<sup>b</sup>, -N(R<sup>a</sup>)C(=O)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)C(=NR<sup>a</sup>)NR<sup>a</sup>R<sup>a</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>R<sup>b</sup>, -N(R<sup>a</sup>)S(=O)<sub>2</sub>NR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylNR<sup>a</sup>R<sup>a</sup>, -NR<sup>a</sup>C<sub>2-6</sub>alkylOR<sup>a</sup> and R<sup>c</sup>.

Claim 38 (new): The compound according to Claim 21, wherein at least one R<sup>6</sup> is selected from C<sub>1-8</sub>alkyl, C<sub>1-4</sub>haloalkyl and halo.

Claim 39 (new): The compound according to Claim 21, wherein at least one R<sup>6</sup> is selected from C<sub>1-4</sub>haloalkyl and halo.

Claim 40 (new): The compound selected from:

2-[4-(2,6-dichlorophenyl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(2,6-dichlorobenzyl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-chloropyridin-4-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole, di-  
trifluoroacetic acid salt,  
4-chloro-6-trifluoromethyl-2-[4-(3-trifluoromethylpyridin-2-yl)piperazin-1-yl]-1H-  
benzoimidazole,  
4-bromo-2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-  
benzoimidazole,  
4-bromo-6-trifluoromethyl-2-[4-(3-trifluoromethylpyridin-2-yl)piperazin-1-yl]-1H-  
benzoimidazole,  
4-chloro-2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-  
benzoimidazole,  
4-bromo-2-[4-(3,5-dichloropyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-  
benzoimidazole,  
4-pyridin-3-yl-6-trifluoromethyl-2-[4-(3-trifluoromethylpyridin-2-yl)piperazin-1-yl]-  
1H-benzoimidazole,  
2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-4-pyridin-3-yl-6-trifluoromethyl-1H-  
benzoimidazole,  
2-[1-(3-chloropyridin-2-yl)piperidin-4-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3,5-dichloropyridin-4-yl)piperazin-1-yl]-6-methyl-1H-benzoimidazole,  
2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-6-methyl-1H-benzoimidazole,  
6-methyl-2-[4-(3-trifluoromethylpyridin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
6-*tert*-butyl-2-[4-(3-chloropyridin-4-yl)piperazin-1-yl]-1H-benzoimidazole,  
6-*tert*-butyl-2-[4-(3,5-dichloropyridin-4-yl)piperazin-1-yl]-1H-benzoimidazole,  
6-*tert*-butyl-2-[4-(3-trifluoromethylpyridin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
6-chloro-2-[4-(3-trifluoromethylpyridin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
6-chloro-2-[4-(3,5-dichloropyridin-4-yl)piperazin-1-yl]-1H-benzoimidazole,  
6-chloro-2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
6-bromo-2-[4-(3-trifluoromethylpyridin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
6-bromo-2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
5,6-dichloro-2-[4-(3-trifluoromethylpyridin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
2-[4-(3-chloropyridin-2-yl)piperazin-1-yl]-5,6-dichloro-1H-benzoimidazole,

6-chloro-2-[4-(3,5-dichloropyridin-4-yl)piperazin-1-yl]-5-methyl-1H-benzoimidazole,  
6-trifluoromethyl-2-[4-(3-trifluoromethylpyridin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
2-[4-(2-chlorophenyl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-methoxypyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
6-Trifluoromethyl-2-[4-(2-trifluoromethylphenylpiperidin-1-yl)-1H-benzoimidazole,  
2-(4-cyano-4-phenylpiperidin-1-yl)-6-trifluoromethyl-1H-benzoimidazole,  
6-trifluoromethyl-2-[4-(3-trifluoromethylphenyl)piperazin-1-yl]-1H-benzoimidazole,  
2-[4-(3,5-dichloropyridin-4-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-(4-phenylpiperazin-1-yl)-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-chlorophenyl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(quinolin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
6-trifluoromethyl-2-[4-(5-trifluoromethylpyridin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
2-[4-(naphthalen-1-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(2,3-dimethylphenyl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(6-methylpyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-chloro-5-trifluoromethylpyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
6-trifluoromethyl-2-[4-(4-trifluoromethylpyrimidin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
2-[4-(pyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
6-trifluoromethyl-2-[4-(4-trifluoromethylbenzyl)piperazin-1-yl]-1H-benzoimidazole,  
6-trifluoromethyl-2-[4-(2-trifluoromethylbenzyl)piperazin-1-yl]-1H-benzoimidazole,  
2-[4-(2,5-dimethylbenzyl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-fluoropyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-chloropyrazin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
6-trifluoromethyl-2-[4-(4-trifluoromethylpyridin-2-yl)piperazin-1-yl]-1H-benzoimidazole,  
2-[4-(3-methylpyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-methylpyridin-2-yl)-[1,4]diazepan-1-yl]-6-trifluoromethyl-1H-benzoimidazole,

2-[4-(3-chloropyridin-2-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-chloropyridin-2-yl)-2,5-dimethylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-chloro-5-hydroxymethylpyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3,5-dichloropyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
(*R*)-2-[4-(3,5-dichloropyridin-2-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
(*S*)-2-[4-(3,5-dichloropyridin-2-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3,5-dichloropyridin-2-yl)-2,5-dimethylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
(*S*)-2-[4-(isoquinolin-1-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
(*S*)-2-[4-(3-methylquinolin-2-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
(*S*)-2-[4-(3-chloro-5-hydroxymethylpyridin-2-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
(*R*)-2-[4-(3-chloro-5-hydroxymethylpyridin-2-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-{4-[3-(pyrrol-1-yl)pyridin-2-yl]piperazin-1-yl}-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(3-iodopyridin-2-yl)piperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
2-[4-(6-trifluoromethyl-1H-benzoimidazol-2-yl)piperazin-1-yl]pyridine-3-sulfonic acid methylamide,  
(*R*)-4-bromo-2-[4-(3,5-dichloropyridin-2-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
(*S*)-4-bromo-2-[4-(3,5-dichloropyridin-2-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
(*S*)-4-bromo-2-[4-(isoquinolin-1-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,



(*S*)-4-bromo-2-[4-(3-chloro-5-hydroxymethylpyridin-2-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
(*R*)-4-bromo-2-[4-(3-chloro-5-hydroxymethylpyridin-2-yl)-2-methylpiperazin-1-yl]-6-trifluoromethyl-1H-benzoimidazole,  
6-trifluoromethyl-4-(3-trifluoromethylphenyl)-2-[4-(3-trifluoromethylpyridin-2-yl)-2-methylpiperazin-1-yl]-1H-benzoimidazole,  
4-(3,4-difluorophenyl)-6-trifluoromethyl-2-[4-(3-trifluoromethylpyridin-2-yl)-2-methylpiperazin-1-yl]-1H-benzoimidazole,  
4-(thiophen-2-yl)-6-trifluoromethyl-2-[4-(3-trifluoromethylpyridin-2-yl)-2-methylpiperazin-1-yl]-1H-benzoimidazole,  
4-(furan-2-yl)-6-trifluoromethyl-2-[4-(3-trifluoromethylpyridin-2-yl)-2-methylpiperazin-1-yl]-1H-benzoimidazole, and  
4-(pyridin-2-yl)-6-trifluoromethyl-2-[4-(3-trifluoromethylpyridin-2-yl)-2-methylpiperazin-1-yl]-1H-benzoimidazole.